





Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; J = New Jersey; O = EPA Office of Water; F = See FAQ; E = Environmental Criteria and Assessment Office; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; R = RBA applied (See User Guide for Arsenic notice); c = cancer; \* = where: n SL < 100X c SL; \*\* = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User Guide); s = Concentration may exceed Csat (See User Guide); SSL values are based on DAF=1

Toxicity and Chemical-specific Information											Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 0.1								
SFO (mg/kg-day) <sup>-1</sup>	K <sub>e</sub> (y <sup>-1</sup> )	IUR (ug/m <sup>3</sup> ) <sup>-1</sup>	K <sub>e</sub> (y <sup>-1</sup> )	RfD <sub>o</sub> (mg/kg-day)	K <sub>e</sub> (y <sup>-1</sup> )	RfC <sub>o</sub> (mg/m <sup>3</sup> )	K <sub>e</sub> (y <sup>-1</sup> )	muta-gen	GIABS	ABS	C <sub>sat</sub> (mg/kg)	PEF (m <sup>3</sup> /kg)	VF (m <sup>3</sup> /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL HQ=0.1 (mg/kg)	Dermal SL HQ=0.1 (mg/kg)	Inhalation SL HQ=0.1 (mg/kg)	Noncarcinogenic SL HI=0.1 (mg/kg)		
				2.0E-03	H									Chloroacetic Acid Chloroacetophenone, 2-	79-11-8 532-27-4					2.3E+02	5.5E+02	1.8E+04	1.6E+02 1.8E+04		
2.0E-01	P			4.0E-03	I									Chloroaniline, p- Chlorobenzene Chlorobenzilate	106-47-8 108-90-7 510-15-6	1.6E+01	3.9E+01		1.1E+01	4.7E+02	1.1E+03	1.4E+02	3.3E+02 1.3E+02 1.6E+03		
1.1E-01	C	3.1E-05	C	2.0E-02	I	5.0E-02	P V				7.6E+02	1.4E+09	6.5E+03	Chlorobenzoic Acid, p- Chlorobenzotrifluoride, 4- Chlorobutane, 1-	74-11-3 98-56-6 109-69-3					3.5E+03	8.3E+03	8.9E+02	2.5E+03 2.5E+02 4.7E+03		
3.1E-02	C	2.3E-05	I	2.0E-02	P	9.8E-02	A V							Chlorodifluoromethane Chloroethanol, 2- Chloroform	75-45-6 107-07-3 67-66-3	1.1E+02		1.4E+00	1.4E+00	2.3E+03		1.1E+02	2.1E+04 2.3E+03 1.0E+02		
2.4E+00	C	6.9E-04	C	3.0E-03	P	1.0E-05	X							Chloromethane Chloromethyl Methyl Ether Chloronitrobenzene, o-	74-87-3 107-30-2 88-73-3	1.4E+00		9.5E-02	8.9E-02	3.5E+02	8.3E+02	6.0E+03	4.6E+01 2.4E+02		
6.3E-03	P			1.0E-03	P	6.0E-04	P							Chloronitrobenzene, p- Chlorophenol, 2- Chloropicrin	100-00-5 95-57-8 76-06-2	5.2E+02	1.2E+03		3.6E+02	1.2E+02	2.8E+02	3.6E+05	8.2E+01 5.8E+02 8.2E-01		
3.1E-03	C	8.9E-07	C	1.5E-02	I									Chloroethanol, 1- Chloroethanol, 2- Chloroethane, 1,1-dichloro-	109-89-7 106-43-4 75-07-6	1.1E+03	2.5E+03	1.9E+07	7.4E+02	1.8E+03	4.1E+03		1.2E+03 2.3E+03 2.3E+03		
2.4E+02	C	6.9E-02	C	2.0E-01	I									Chloroethane, 1,1-dichloro- Chloroethane, 1,2-dichloro- Chloroethane, 1,1,1-trichloro-	101-21-3 2921-88-2 5598-13-0	1.4E-02	3.2E-02	2.4E+02	9.6E-03	2.3E+04	5.5E+04	2.8E+02	1.6E+04 8.2E+01		
5.0E-01	J	8.4E-02	S	1.5E+00	I									Chloroethane, 1,1,1-trichloro- Chloroethane, 1,2-dichloro- Chloroethane, 1,2-dichloro-	100-00-5 95-57-8 76-06-2	6.5E+00		2.0E+02	6.3E+00	1.2E+02	2.8E+02	6.0E+04	8.2E+01 5.8E+02 8.2E-01		
9.0E-03	P	3.0E-04	P	3.0E-03	I	1.0E-04	I	M						Chloroethane, 1,1,1-trichloro- Chloroethane, 1,2-dichloro- Chloroethane, 1,2-dichloro-	100-00-5 95-57-8 76-06-2			1.9E+03	1.9E+03	3.5E+01		3.6E+03	3.5E+01		
6.2E-04	I			4.0E-02	H									Cobalt Coke Oven Emissions Copper	7440-48-4 8007-45-2 7440-50-8					4.7E+03			4.7E+03		
1.9E+00	H			5.0E-02	I	6.0E-01	C							Cresol, m- Cresol, o- Cresol, p- Cresol, p-chloro- Cresols Crotonaldehyde, trans-	108-39-4 95-48-7 106-44-5 59-50-7 1319-77-3 123-73-9	1.1E+00			1.7E+00			5.8E+03 5.8E+03 1.2E+04 1.2E+04 1.2E+02	1.4E+04 1.4E+04 2.8E+04 2.8E+04	3.6E+08 3.6E+08 3.6E+08	4.1E+03 4.1E+03 8.2E+03 8.2E+03 1.2E+02
2.2E-01	C	6.3E-05	C	1.0E-01	I	4.0E-01	I V							Cumene Cupferron Cyanazine	98-82-8 135-20-6 21725-46-2	1.5E+01	3.5E+01	2.6E+05	1.0E+01	2.3E+02	5.5E+02		1.1E+03	9.9E+02 1.6E+02	
8.4E-01	H			2.0E-03	H									<b>Cyanides</b> -Calcium Cyanide -Copper Cyanide	592-01-8 544-92-3					1.2E+02	5.8E+02		1.2E+02 5.8E+02		
				6.0E-04	I	8.0E-04	S V							-Cyanide (CN-) -Cyanogen -Cyanogen Bromide	57-12-5 460-19-5 506-68-3					7.0E+01		1.2E+00	1.2E+00 1.2E+02 1.1E+04		
				5.0E-02	I									-Cyanogen Chloride -Hydrogen Cyanide -Potassium Cyanide	506-77-4 74-90-8 151-50-8					5.8E+03		1.8E+01	5.8E+03 1.5E+01 2.3E+02		
				5.0E-03	I	4.0E-02								-Potassium Silver Cyanide -Silver Cyanide -Sodium Cyanide	506-61-6 506-64-9 143-33-9					5.8E+02			5.8E+02 1.2E+04 1.2E+02		
				2.0E-04	P									-Thiocyanates -Thiocyanic Acid -Zinc Cyanide	NA 463-56-9 557-21-1					2.3E+01			2.3E+01 2.3E+01 5.8E+03		
2.3E-02	H			6.0E+00	I V									Cyclohexane Cyclohexane, 1,2,3,4,5-pentabromo-6-chloro- Cyclohexanone	110-82-7 87-84-3 108-94-1	1.4E+02	3.4E+02		1.0E+02			2.7E+03	2.7E+03		
				5.0E+00	I	7.0E-01	P V							Cyclohexene Cyclohexylamine Cylhalothrin/karate	110-83-8 108-91-8 68085-85-8					5.8E+02		6.4E+02	3.1E+02 2.3E+04 4.1E+02		
2.4E-01	I	6.9E-05	C	1.0E-02	I									Cypermethrin Cyromazine DDD	52315-07-8 66215-27-8 72-54-8	1.4E+01	3.2E+01	2.4E+05	9.6E+00	1.2E+03	2.8E+03	2.1E+03	8.2E+02 6.2E+02		





Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; J = New Jersey; O = EPA Office of Water; F = See FAQ; E = Environmental Criteria and Assessment Office; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; R = RBA applied (See User Guide for Arsenic notice); c = cancer; \* = where: n SL < 100X c SL; \*\* = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User Guide); s = Concentration may exceed Csat (See User Guide); SSL values are based on DAF=1

Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 0.1			
SFO (mg/kg-day) <sup>-1</sup>	K <sub>e</sub> (y <sup>-1</sup> )	IUR (ug/m <sup>3</sup> -y) <sup>-1</sup>	K <sub>e</sub> (y <sup>-1</sup> )	RfD <sub>o</sub> (mg/kg-day)	RfC <sub>i</sub> (mg/m <sup>3</sup> )	K <sub>e</sub> (y <sup>-1</sup> )	muta- gen	GIABS	ABS	C <sub>sat</sub> (mg/kg)	PEF (m <sup>3</sup> /kg)	VF (m <sup>3</sup> /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL HQ=0.1 (mg/kg)	Dermal SL HQ=0.1 (mg/kg)	Inhalation SL HQ=0.1 (mg/kg)	Noncarcinogenic SL HI=0.1 (mg/kg)
				8.0E-02 2.0E-02	I				1.0E+00 1.0E-01	1.0E-01 1.0E-01	1.4E+09 1.4E+09		Fluridone Flurprimidol	59756-60-4 56425-91-3					9.3E+03 2.3E+03	2.2E+04 5.5E+03		6.6E+03 1.6E+03
3.5E-03	I			6.0E-02 1.0E-02 1.0E-01	I				1.0E+00 1.0E+00 1.0E+00	1.0E-01 1.0E-01 1.0E-01	1.4E+09 1.4E+09 1.4E+09		Flutolanil Fluvalinate Folpet	66332-96-5 69409-94-5 133-07-3	9.3E+02	2.2E+03		6.6E+02	7.0E+03 1.2E+03 1.2E+04	1.7E+04 2.8E+03 2.8E+04		4.9E+03 8.2E+02 8.2E+03
1.9E-01	I			2.0E-03 2.0E-01	I				1.0E+00 1.0E+00	1.0E-01 1.0E-01	1.4E+09 1.4E+09		Fomesafen Fonofos Formaldehyde	72178-02-0 944-22-9 50-00-0	1.7E+01	4.1E+01		1.2E+01	2.3E+02 2.3E+04	5.5E+02		1.6E+02 3.3E+02
		1.3E-05	I	9.0E-01 3.0E+00	P	3.0E-04	X	V	1.0E+00		1.1E+05 1.4E+09	9.3E+04	Formic Acid Fosetyl-AL Furans	64-18-6 39148-24-8				7.3E+01	7.3E+01		1.2E+01 1.2E+05	1.2E+01 2.5E+05
				1.0E-03 1.0E-03 9.0E-01	X			V	1.0E+00 1.0E+00 1.0E+00	3.0E-02 3.0E-02 3.0E-02	1.4E+09 6.2E+03 1.4E+09	2.0E+05 2.6E+03 1.2E+04	Dibenzofuran Furan Tetrahydrofuran	132-64-9 110-00-9 109-99-9					1.2E+02 1.2E+02 1.1E+05	9.2E+02 9.2E+02 8.3E+05	1.1E+04	1.0E+02 1.0E+02 9.6E+03
3.8E+00	H			3.0E-03	I	5.0E-02	H	V	1.0E+00		1.0E+04	4.9E+04	Furazolidone Furfural Furium	67-45-8 98-01-1 531-82-8	8.6E-01	2.0E+00		6.0E-01	3.5E+02		1.1E+03	2.6E+02
1.5E+00	C	4.3E-04	C						1.0E+00	1.0E-01	1.4E+09		Furmecycloz Glufosinate, Ammonium Glutaraldehyde	60568-05-0 77182-82-2 111-30-8	1.1E+02	2.6E+02	1.9E+06	7.7E+01	4.7E+01	1.1E+02		3.3E+01 4.8E+04
				4.0E-04 1.0E-01 3.0E-03	I	1.0E-03	H	V	1.0E+00		1.1E+05 1.4E+09	7.3E+04	Glycidyl Glyphosate Goal	765-34-4 1071-83-6 42874-03-3					4.7E+01 1.2E+04 3.5E+02	3.2E+01		1.9E+01 8.2E+03 2.5E+02
				1.0E-02 2.0E-02 3.0E-03	X			V	1.0E+00		1.4E+09	1.5E+05	Guandine Guandine Chloride Guthion	113-00-8 50-01-1 98-50-0					1.2E+03 2.3E+03 3.5E+02	5.5E+03 8.3E+02	6.0E+06	1.2E+03 1.6E+03 2.5E+02
4.5E+00	I	1.3E-03	I	5.0E-05 1.3E-02 5.0E-04	I	1.0E-01	H	V	1.0E+00		1.4E+09		Haloxypol, Methyl Harmony Heptachlor	69806-40-2 79277-27-3 76-44-8	7.3E-01		4.5E+00	6.3E-01	5.8E+00 1.5E+03 5.8E+01	1.4E+01 3.6E+03		4.1E+00 1.1E+03 5.8E+01
9.1E+00	I	2.6E-03	I	1.3E-05 2.0E-03 2.0E-04	I			V	1.0E+00		1.4E+09	8.4E+05	Heptachlor Epoxide Hexabromobenzene Hexabromodiphenyl ether, 2,2',4,4',5,5'- (BDE-153)	1024-57-3 87-82-1 68631-49-2	3.6E-01		4.0E+00	3.3E-01	1.5E+00 2.3E+02 2.3E+01		5.5E+01	1.5E+00 2.3E+02 1.6E+01
1.6E+00	I	4.6E-04	I	8.0E-04	I			V	1.0E+00		1.4E+09	6.8E+04	Hexachlorobenzene Hexachlorobutadiene Hexachlorocyclohexane, Alpha-	118-74-1 87-68-3 319-84-6	2.0E-00 4.2E-01 5.2E-01		1.8E+00 6.0E+00 1.2E+00	9.6E-01 5.3E+00 3.6E-01	9.3E+01 1.2E+02 9.3E+02	2.2E+03		9.3E+01 1.2E+02 6.6E+02
1.8E+00	I	5.3E-04	I						1.0E+00	1.0E-01	1.4E+09		Hexachlorocyclohexane, Beta- Hexachlorocyclohexane, Gamma- (Lindane) Hexachlorocyclohexane, Technical	319-85-7 58-89-9 508-73-1	1.8E-00 3.0E-00 1.8E-00	4.3E+00 1.8E+01 4.3E+00	3.1E+04 5.4E+04 3.3E+04	1.3E+00 2.5E+00 1.3E+00	3.5E+01	2.1E+02		3.0E+01
4.0E-02	I	1.1E-05	C	6.0E-03 7.0E-04 3.0E-04	I	2.0E-04	I	V	1.0E+00		1.6E+01 1.4E+09	8.5E+03	Hexachlorocyclopentadiene Hexachloroethane Hexachlorophene	77-47-4 67-72-1 70-30-4	8.2E-01		8.9E+00	8.0E+00	7.0E+02 8.2E+01 3.5E+01	7.5E-01 1.1E+02		7.5E-01 4.6E+01 2.5E+01
1.1E-01	I			3.0E-03 4.0E-04	I	1.0E-05	I	V	1.0E+00		5.2E+03 1.4E+09	3.0E+05	Hexahydro-1,3,5-trinitro-1,3,5-triazine (RDX) Hexamethylene Diisocyanate, 1,6- Hexamethylphosphoramide	121-82-4 822-06-0 680-31-9	3.0E-01	4.7E+02		2.8E+01	3.5E+02 4.7E+01	5.5E+03 1.1E+02		3.3E+02 1.3E+00 3.3E+01
				6.0E-02 2.0E+00 5.0E-03	H	7.0E-01	I	V	1.0E+00		1.4E+02 1.4E+09	8.3E+02	Hexane, N- Hexanedioic Acid Hexanone, 2-	110-54-3 124-04-9 591-78-6					7.0E+03 2.3E+05 5.8E+02	5.5E+05 1.7E+02		2.5E+02 1.6E+05 1.3E+02
3.0E+00	I	4.9E-03	I						1.0E+00	1.0E-01	1.4E+09		Hexazinone Hydrazine Hydrazine Sulfate	51235-04-2 302-01-2 10034-93-2	1.1E+00 1.1E+00		3.4E+03 3.4E+03	1.1E+00 1.1E+00	3.9E+03	9.1E+03		2.7E+03 1.8E+04
3.0E+00	I	4.9E-03	I						1.0E+00	1.0E-01	1.4E+09		Hydrogen Chloride Hydrogen Fluoride Hydrogen Sulfide	7647-01-0 7664-39-3 7783-06-4				4.7E+03			1.2E+07 8.3E+06 1.2E+06	1.2E+07 4.7E+03 1.2E+06
6.0E-02	P			4.0E-02 1.3E-02 2.5E-01	P			V	1.0E+00	1.0E-01	1.4E+09		Hydroquinone Imazail Imazaquin	123-31-9 35554-44-0 81335-37-7	5.5E+01	1.3E+02		3.8E+01	4.7E+03 1.5E+03 2.9E+04	1.1E+04 3.6E+03 6.9E+04		3.3E+03 1.1E+03 2.1E+04
				1.0E-02 4.0E-02 7.0E-01	A				1.0E+00		1.4E+09		Iodine Iprodione Iron	7553-56-2 36734-19-7 7439-89-6					1.2E+03 4.7E+03 8.2E+04	1.1E+04		1.2E+03 3.3E+03 8.2E+04
9.5E-04	I			3.0E-01 2.0E-01 1.5E-02	I	2.0E+00	C	V	1.0E+00	1.0E-01	1.4E+09	2.8E+04	Isobutyl Alcohol Isophorone Isopropalin	78-83-1 78-59-1 33820-53-0	3.4E+03	8.1E+03		2.4E+03	3.5E+04 2.3E+04 1.8E+03	5.5E+04	1.2E+09	3.5E+04 1.6E+04 1.8E+03
				2.0E+00 1.0E-01 5.0E-02	P	2.0E-01	P	V	1.0E+00		1.1E+05 1.4E+09	2.8E+04	Isopropanol Isopropyl Methyl Phosphonic Acid Isoxaben	67-63-0 1832-54-8 82558-50-7					2.3E+05 1.2E+04 5.8E+03	2.4E+03 2.8E+04 1.4E+04		2.4E+03 8.2E+03 4.1E+03













Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; J = New Jersey; O = EPA Office of Water; F = See FAQ; E = Environmental Criteria and Assessment Office; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; R = RBA applied (See User Guide for Arsenic notice); c = cancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User Guide); s = Concentration may exceed Csat (See User Guide); SSL values are based on DAF=1																							
Toxicity and Chemical-specific Information										Contaminant			Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 0.1						
SFO	ke	IUR	ke	RfD	ke	RfC	ke	muta-	GIABS	ABS	Csat	PEF	VF	Analyte	CAS No.	Ingestion SL TR=1.0E-6	Dermal SL TR=1.0E-6	Inhalation SL TR=1.0E-6	Carcinogenic SL TR=1.0E-6	Ingestion SL HQ=0.1	Dermal SL HQ=0.1	Inhalation SL HQ=0.1	Noncarcinogenic SL HI=0.1
(mg/kg-day) <sup>-1</sup>	y	(ug/m <sup>3</sup> ) <sup>-1</sup>	y	(mg/kg-day)	y	(mg/m <sup>3</sup> )	y	gen	(mg/kg)	(mg/kg)	(mg/kg)	(m <sup>3</sup> /kg)	(m <sup>3</sup> /kg)			(mg/kg)	(mg/kg)	(mg/kg)	(mg/kg)	(mg/kg)	(mg/kg)	(mg/kg)	(mg/kg)
				2.0E-03 7.0E-06	P				1.0E+00 1.0E+00	6.5E-04		1.4E+09 1.4E+09		Tetryl (Trinitrophenylmethylintramine) Thallium (I) Nitrate	479-45-8 10102-45-1					2.3E+02 8.2E-01		8.5E+04	2.3E+02 8.2E-01
				1.0E-05 6.0E-06 2.0E-05	X X X				1.0E+00 1.0E+00 1.0E+00	1.0E-01 1.0E-01		1.4E+09 1.4E+09 1.4E+09		Thallium (Soluble Salts) Thallium Acetate Thallium Carbonate	7440-28-0 563-68-8 6533-73-9					1.2E+00 7.0E-01 2.3E+00	1.7E+00 5.5E+00		1.2E+00 4.9E-01 1.6E+00
				6.0E-06 2.0E-05 1.0E-02	X X I				1.0E+00 1.0E+00 1.0E+00	1.0E-01 1.0E-01		1.4E+09 1.4E+09 1.4E+09		Thallium Chloride Thallium Sulfate Thiocarb	7791-12-0 7446-18-6 28249-77-6					7.0E-01 2.3E+00 1.2E+03	2.6E+05 2.8E+03		7.0E-01 2.3E+00 8.2E+02
				7.0E-02 3.0E-04 8.0E-02	X H I				1.0E+00 1.0E+00 1.0E+00	7.5E-03 1.0E-01		1.4E+09 1.4E+09 1.4E+09		Thiodiglycol Thiofanox Thiophanate, Methyl	111-48-8 39196-18-4 23564-05-8					8.2E+03 3.5E+01 9.3E+03	2.6E+05 8.3E+01 2.2E+04		7.9E+03 2.5E+01 6.6E+03
				5.0E-03 6.0E-01	I H				1.0E+00 1.0E+00 1.0E+00	1.0E-01		1.4E+09 1.4E+09 1.4E+09		Thiram Tin Titanium Tetrachloride	137-26-8 7440-31-5 7550-45-0					5.8E+02 7.0E+04	1.4E+03	6.0E+04	4.1E+02 7.0E+04 6.0E+04
				8.0E-02 1.8E-01 3.0E-02	I X P				1.0E+00 1.0E+00 1.0E+00	8.2E+02 1.0E-01		4.3E+03 1.4E+09 1.4E+09		Toluene Toluene-2,5-diamine Toluidine, p-	108-88-3 95-70-5 106-49-0	1.8E+01 1.1E+02	4.3E+01 2.6E+02		1.3E+01 7.7E+01	9.3E+03 2.3E+01 4.7E+02	5.5E+01 1.1E+03		4.7E+03 1.6E+01 3.3E+02
				3.0E+00 1.0E-02 4.0E-02	P V P				1.0E+00 1.0E+00 1.0E+00	3.4E-01 1.4E+02 6.9E+00		1.4E+09 8.3E+02 1.0E+03	1.1E+03	Total Petroleum Hydrocarbons (Aliphatic High) Total Petroleum Hydrocarbons (Aliphatic Low) Total Petroleum Hydrocarbons (Aliphatic Medium)	NA NA NA					3.5E+05 1.2E+03		2.2E+02 4.6E+01	3.5E+05 2.2E+02 4.4E+01
				4.0E-02 4.0E-03 4.0E-03	P P P				1.0E+00 1.0E+00 1.0E+00	1.0E-01 3.0E-02 3.0E-03		1.4E+09 1.8E+03 5.2E+04	1.4E+09 3.5E+03	Total Petroleum Hydrocarbons (Aromatic High) Total Petroleum Hydrocarbons (Aromatic Low) Total Petroleum Hydrocarbons (Aromatic Medium)	NA NA NA					4.7E+03 4.7E+02 4.7E+02	1.1E+04	4.6E+01	3.3E+03 4.2E+01 6.0E+01
				1.1E+00 7.5E-03 3.0E-04	I I A				1.0E+00 1.0E+00 1.0E+00	1.0E-01 1.0E-01		1.4E+09 1.4E+09 1.4E+09	3.4E+03	Toxaphene Tralomethrin Tri-n-butyltin	8001-35-2 66841-25-6 688-73-3	3.0E+00 7.0E+00	5.2E+04	2.1E+00	8.8E+02 3.5E+01	2.1E+03		6.2E+02 3.5E+01	
				8.0E+01 1.3E-02 1.0E-02	X I I				1.0E+00 1.0E+00 1.0E+00	1.0E-01 1.0E-01		1.4E+09 3.6E+05	4.5E+04	Triacetin Triallate Trisulfuron	102-76-1 2303-17-5 82499-40-5					9.3E+06 1.5E+03 1.2E+03	2.2E+07		6.6E+06 1.5E+03 8.2E+02
				5.0E-03 1.0E-02 3.0E-04	I P P				1.0E+00 1.0E+00 1.0E+00	1.0E-01 1.0E-01		1.4E+09 1.4E+09	4.5E+04	Tribromobenzene, 1,2,4- Tributyl Phosphate Tributyltin Compounds	815-64-3 126-73-8 NA	3.6E+02	8.6E+02	2.6E+02	5.8E+02 1.2E+03 3.5E+01	2.8E+03	8.3E+01	5.8E+02 8.2E+02 2.5E+01	
				3.0E-04 3.0E+01 7.0E-02	I I I				1.0E+00 1.0E+00 1.0E+00	1.0E-01 3.0E+01	9.1E+02	1.4E+09 1.3E+03	1.3E+03	Tributyltin Oxide Trichloro-1,2,2-trifluoroethane, 1,1,2- Trichloroacetic Acid	56-35-9 76-13-1 76-03-9	4.7E+01	1.1E+02	3.3E+01	3.5E+01 3.5E+06 2.3E+03	8.3E+01	1.7E+04	2.5E+01 1.7E+04 1.6E+03	
				2.9E-02 7.0E-03 8.0E-04	H X X				1.0E+00 1.0E+00 1.0E+00	1.0E-01 1.0E-01		1.4E+09 1.4E+09 3.2E+04	3.2E+04	Trichloroamine HCl, 2,4,6- Trichloroaniline, 2,4,6- Trichlorobenzene, 1,2,3-	33663-50-2 634-93-5 87-61-6	4.1E+02 1.7E+02	2.7E+02 1.1E+03	7.9E+01 3.3E+02	3.5E+00 9.3E+01	8.3E+00		2.5E+00 9.3E+01	
				2.9E-02 2.0E+00 5.7E-02	P I I				1.0E+00 1.0E+00 1.0E+00	2.0E-03 5.0E+00 2.0E-04		3.0E+04 1.7E+03 7.2E+03	3.0E+04	Trichlorobenzene, 1,2,4- Trichloropropane, 1,1,1- Trichloroethane, 1,1,2-	120-82-1 71-55-6 79-00-5	1.1E+02 5.7E+01	5.5E+00	5.0E+00	1.2E+03 2.3E+05 4.7E+02	2.6E+01 3.6E+03 6.3E-01	2.6E+01 3.6E+03 6.3E-01		
				4.6E-02 3.0E-01 1.0E-01	I I I				1.0E+00 1.0E+00 1.0E+00	2.0E-03 7.0E-01		2.2E+03 1.0E+03	2.2E+03	Trichloroethylene Trichlorofluoromethane Trichlorophenol, 2,4,5-	79-01-8 75-69-4 95-95-4	7.1E+01	6.6E+00	6.0E+00	5.8E+01 3.5E+04 1.2E+04	1.9E+00 3.2E+02		1.9E+00 3.1E+02 8.2E+03	
				1.1E-02 1.0E-02 8.0E-03	I I I				1.0E+00 1.0E+00 1.0E+00	1.0E-01 1.0E-01		1.4E+09 1.4E+09 1.4E+09		Trichlorophenol, 2,4,6- Trichlorophenoxyacetic Acid, 2,4,5- Trichlorophenoxypropionic acid, -2,4,5	88-06-2 93-76-5 93-72-1	3.0E+02	7.0E+02	5.4E+06	2.1E+02	1.2E+02 1.2E+03 9.3E+02	2.8E+02 2.8E+03		8.2E+01 8.2E+02 6.6E+02
				3.0E+01	I				1.0E+00	1.3E+03		1.4E+09	1.5E+04	Trichloropropane, 1,1,2- Trichloropropane, 1,2,3- Trichloropropene, 1,2,3-	598-77-6 96-18-4 96-19-5	1.1E-01		1.1E-01	5.8E+02 4.7E+02 3.5E+02	2.1E+00 3.1E-01		5.8E+02 2.1E+00 3.1E-01	
				2.0E-02 3.0E-03	A I				1.0E+00 1.0E+00	1.0E-01		1.4E+09 1.4E+09	1.6E+04	Tricresyl Phosphate (TCP) Triethyldiamine	1330-78-5 58138-08-2 121-44-8					2.3E+03 3.5E+02	5.5E+03 8.3E+02	4.8E+01	1.8E+03 2.5E+02 4.8E+01
				7.7E-03 2.0E-02	I P				1.0E+00 1.0E+00	1.0E-01		1.4E+09 5.1E+05	5.1E+05	Triethylene Glycol Trifluralin Trimethyl Phosphate	112-27-6 1582-09-8 512-56-1	4.2E+02 1.6E+02	3.9E+02		4.2E+02 1.1E+02	8.8E+02 1.2E+03	2.8E+03		1.6E+05 8.8E+02 8.2E+02
				5.0E-03 7.0E-03 1.0E-02	P P X				1.0E+00 1.0E+00 1.0E+00	2.9E+02 2.2E+02 1.8E+02		9.4E+03 7.8E+03 6.6E+03	9.4E+03	Trimethylbenzene, 1,2,3- Trimethylbenzene, 1,2,4- Trimethylbenzene, 1,3,5-	525-73-8 95-63-6 108-67-8						2.1E+01 2.4E+01		2.1E+01 2.4E+01 1.2E+03
				3.0E-02	I				1.0E+00	1.9E-02		1.4E+09		Trinitrobenzene, 1,3,5- Trinitrotoluene, 2,4,6- Triphenylphosphine Oxide	99-35-4 118-96-7 791-28-6	1.1E+02	8.0E+02	9.6E+01	3.5E+03 5.8E+01 2.3E+03	4.4E+04 4.3E+02 5.5E+03		3.2E+03 5.1E+01 1.6E+03	
				2.0E-02 1.0E-02	A X				1.0E+00 1.0E+00	1.0E-01		1.4E+09	1.4E+09	Tris(1,3-Dichloro-2-propyl) Phosphate Tris(1-chloro-2-propyl)phosphate Tris(2,3-dibromopropyl)phosphate	13674-87-8 13674-84-5 126-72-7	1.4E+00		1.7E+01	1.3E+00	2.3E+03 1.2E+03	5.5E+03 2.8E+03		1.6E+03 8.2E+02

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; J = New Jersey; O = EPA Office of Water; F = See FAQ; E = Environmental Criteria and Assessment Office; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; R = RBA applied (See User Guide for Arsenic notice); c = cancer; \* = where n SL < 100X c SL; \*\* = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User Guide); s = Concentration may exceed Csat (See User Guide); SSL values are based on DAF=1

Toxicity and Chemical-specific Information											Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 0.1							
SFO (mg/kg-day) <sup>-1</sup>	Ke (y)	IUR (ug/m <sup>3</sup> ) <sup>-1</sup>	Ke (y)	RfD <sub>o</sub> (mg/kg-day)	Ke (y)	RfC <sub>i</sub> (mg/m <sup>3</sup> )	Ke (y)	muta- gen	GIABS	ABS	C <sub>sat</sub> (mg/kg)	PEF (m <sup>3</sup> /kg)	VF (m <sup>3</sup> /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL HQ=0.1 (mg/kg)	Dermal SL HQ=0.1 (mg/kg)	Inhalation SL HQ=0.1 (mg/kg)	Noncarcinogenic SL HI=0.1 (mg/kg)	
2.0E-02	P			7.0E-03	P									Tris(2-chloroethyl)phosphate	115-96-8	1.6E+02	3.9E+02		1.1E+02	8.2E+02	1.9E+03		5.7E+02	
3.2E-03	P			1.0E-01	P									Tris(2-ethylhexyl)phosphate	78-42-2	1.0E+03	2.4E+03		7.2E+02	1.2E+04	2.8E+04		8.2E+03	
				3.0E-03	I	4.0E-05	A							Uranium (Soluble Salts)	NA				3.5E+02		2.4E+04		3.5E+02	
1.0E+00	C	2.9E-04	C					M	1.0E+00	1.0E-01				Urethane	51-79-6	3.3E+00	7.7E+00	5.7E+04	2.3E+00	1.1E+03			8.4E+02	
		8.3E-03	P	9.0E-03	I	7.0E-06	P		2.6E-02					Vanadium Pentoxide	1314-62-1			2.0E+03	2.0E+03	5.9E+02		4.2E+03	8.4E+02	
				5.0E-03	S	1.0E-04	A		2.6E-02					Vanadium and Compounds	7440-62-2					5.9E+02		6.0E+04	5.8E+02	
				1.0E-03	I		V		1.0E+00				1.2E+05	Vermolate	1929-77-7					1.2E+02			1.2E+02	
				2.5E-02	I				1.0E+00	1.0E-01				Vinclozolin	50471-44-8					2.9E+03	6.9E+03		2.1E+03	
				1.0E+00	H	2.0E-01	I	V	1.0E+00		2.8E+03	1.4E+09	4.4E+03	Vinyl Acetate	108-05-4					1.2E+05			3.9E+02	3.8E+02
		3.2E-05	H			3.0E-03	I	V	1.0E+00		3.4E+03	1.4E+09	1.4E+03	Vinyl Bromide	593-60-2			5.2E-01	5.2E-01			1.8E+00	1.8E+00	
7.2E-01	I	4.4E-06	I	3.0E-03	I	1.0E-01	I	V	M	1.0E+00	3.9E+03	1.4E+09	9.6E+02	Vinyl Chloride	75-01-4	4.5E+00		2.7E+00	1.7E+00	3.5E+02		4.2E+01	3.7E+01	
				3.0E-04	I				1.0E+00	1.0E-01				Warfarin	81-81-2					3.5E+01	8.3E+01		2.5E+01	
				2.0E-01	S	1.0E-01	S	V	1.0E+00		3.9E+02	1.4E+09	5.6E+03	Xylene, p-	106-42-3					2.3E+04		2.4E+02	2.4E+02	
				2.0E-01	S	1.0E-01	S	V	1.0E+00		3.9E+02	1.4E+09	5.5E+03	Xylene, m-	108-38-3					2.3E+04		2.4E+02	2.4E+02	
				2.0E-01	S	1.0E-01	S	V	1.0E+00		4.3E+02	1.4E+09	6.5E+03	Xylene, o-	95-47-6					2.3E+04		2.8E+02	2.8E+02	
				2.0E-01	I	1.0E-01	I	V	1.0E+00		2.6E+02	1.4E+09	6.5E+03	Xylenes	1330-20-7					2.3E+04		2.8E+02	2.8E+02	
				3.0E-04	I				1.0E+00			1.4E+09		Zinc Phosphide	1314-84-7					3.5E+01			3.5E+01	
				3.0E-01	I				1.0E+00			1.4E+09		Zinc and Compounds	7440-66-6					3.5E+04			3.5E+04	
				5.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Zincb	12122-67-7					5.8E+03	1.4E+04		4.1E+03	
				8.0E-05	X				1.0E+00			1.4E+09		Zirconium	7440-67-7					9.3E+00			9.3E+00	